

Classical Unsupervised Machine Learning

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Different Types of Machine Learning

	Supervised	Unsupervised
Classical	Decision Trees, SVMs, shallow neural networks	Clustering, Dimensionality Reduction
Deep	Multi-layer ANNs/ CNNs, ResNets	GANs, VAEs, INNs

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- Unsupervised = allowing the computer to identify the important features in your data.
- This is beneficial for identifying patterns or important quantities in your data.
- Prior knowledge of the dataset isn't required!

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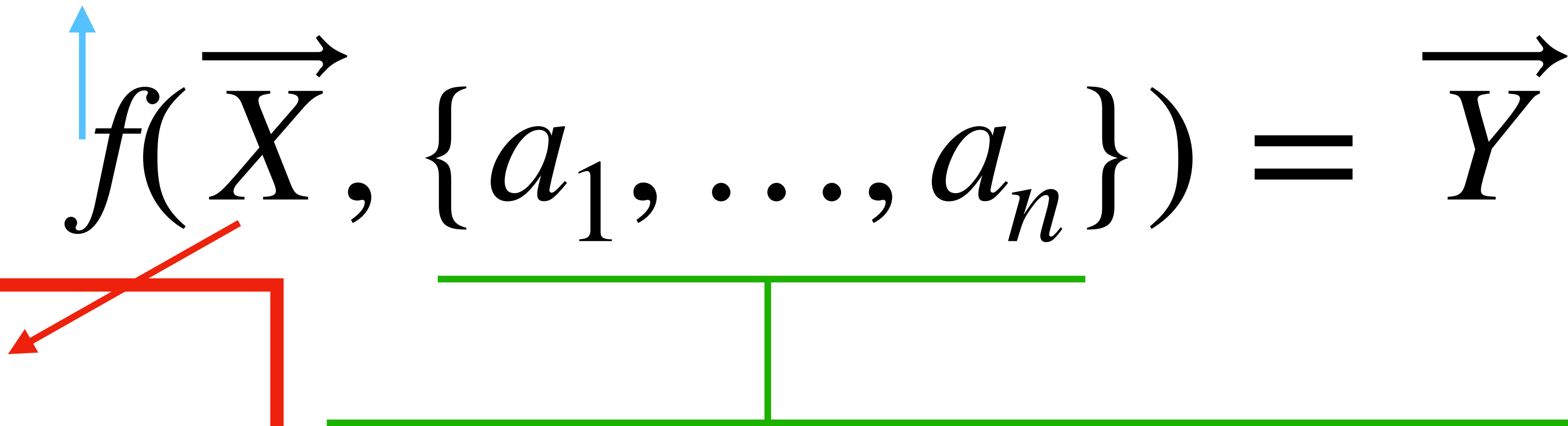
Hyperparameters:

- Tuning parameters for the algorithm.
- The supervised bit of unsupervised learning

How does unsupervised learning work?

Internal Choices of the computer:

- How the computer manipulates the input and the hyperparameters

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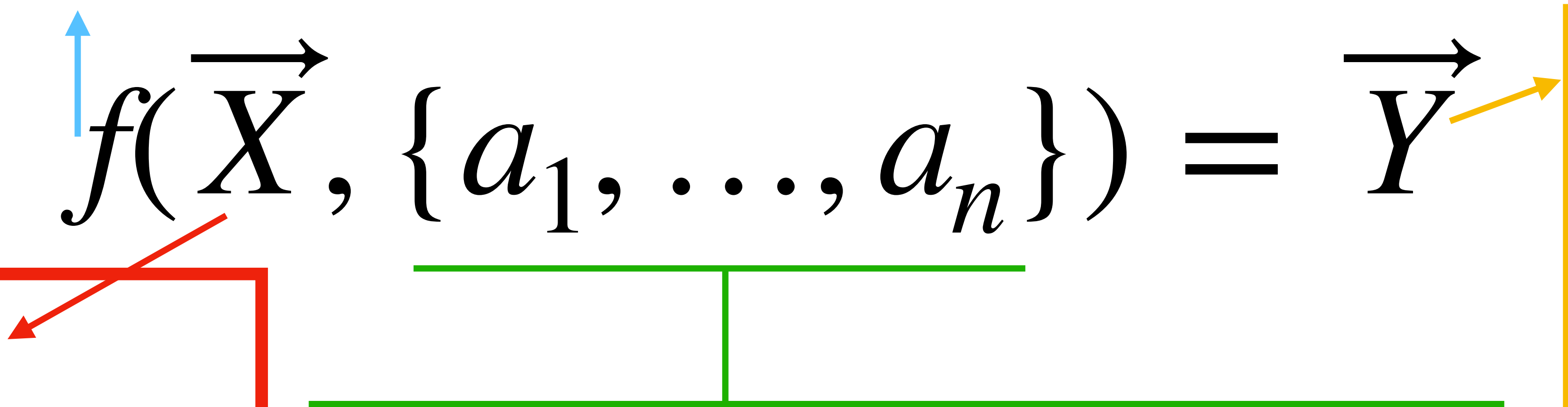
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Output:

- The output of the algorithm

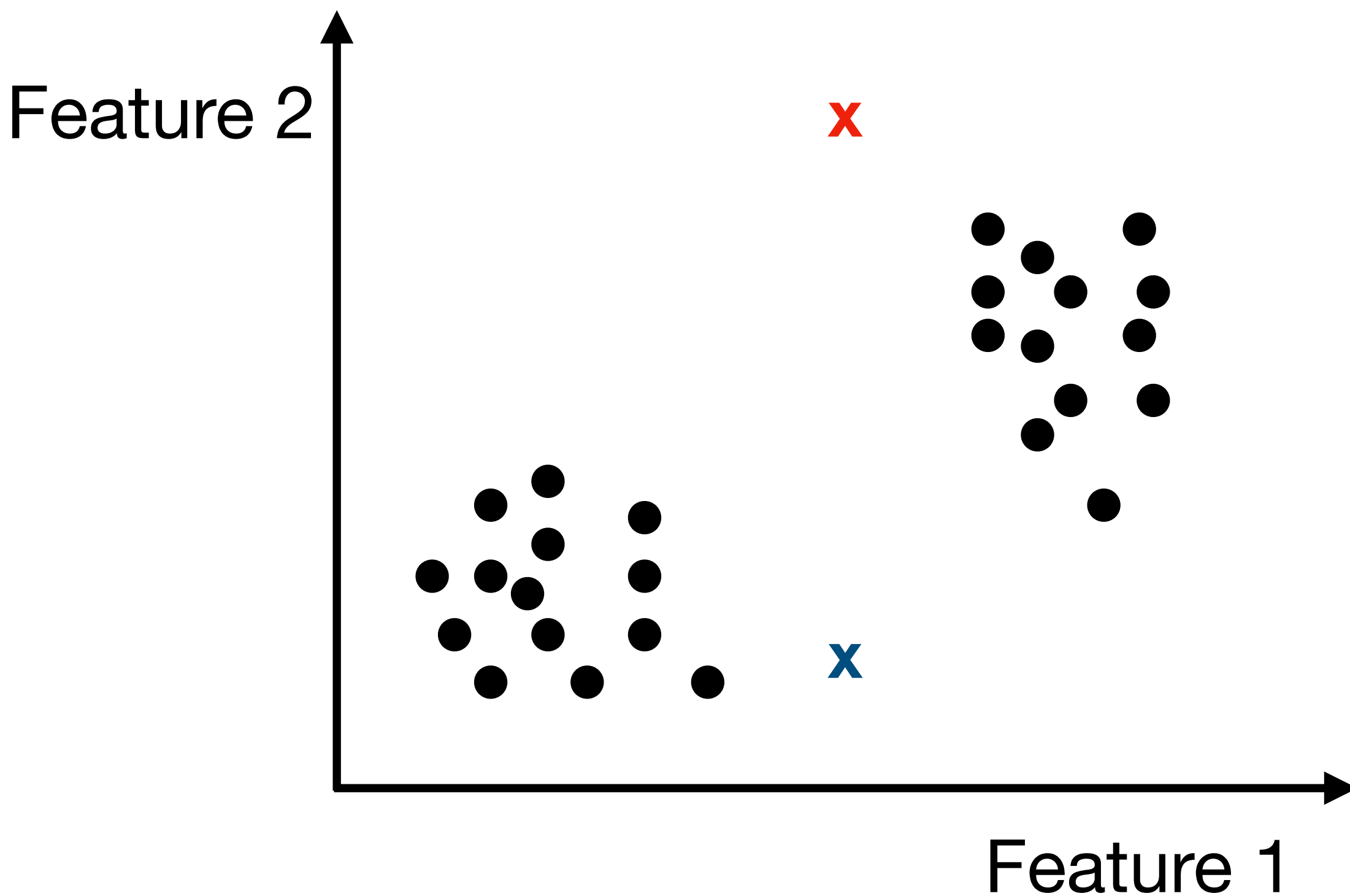
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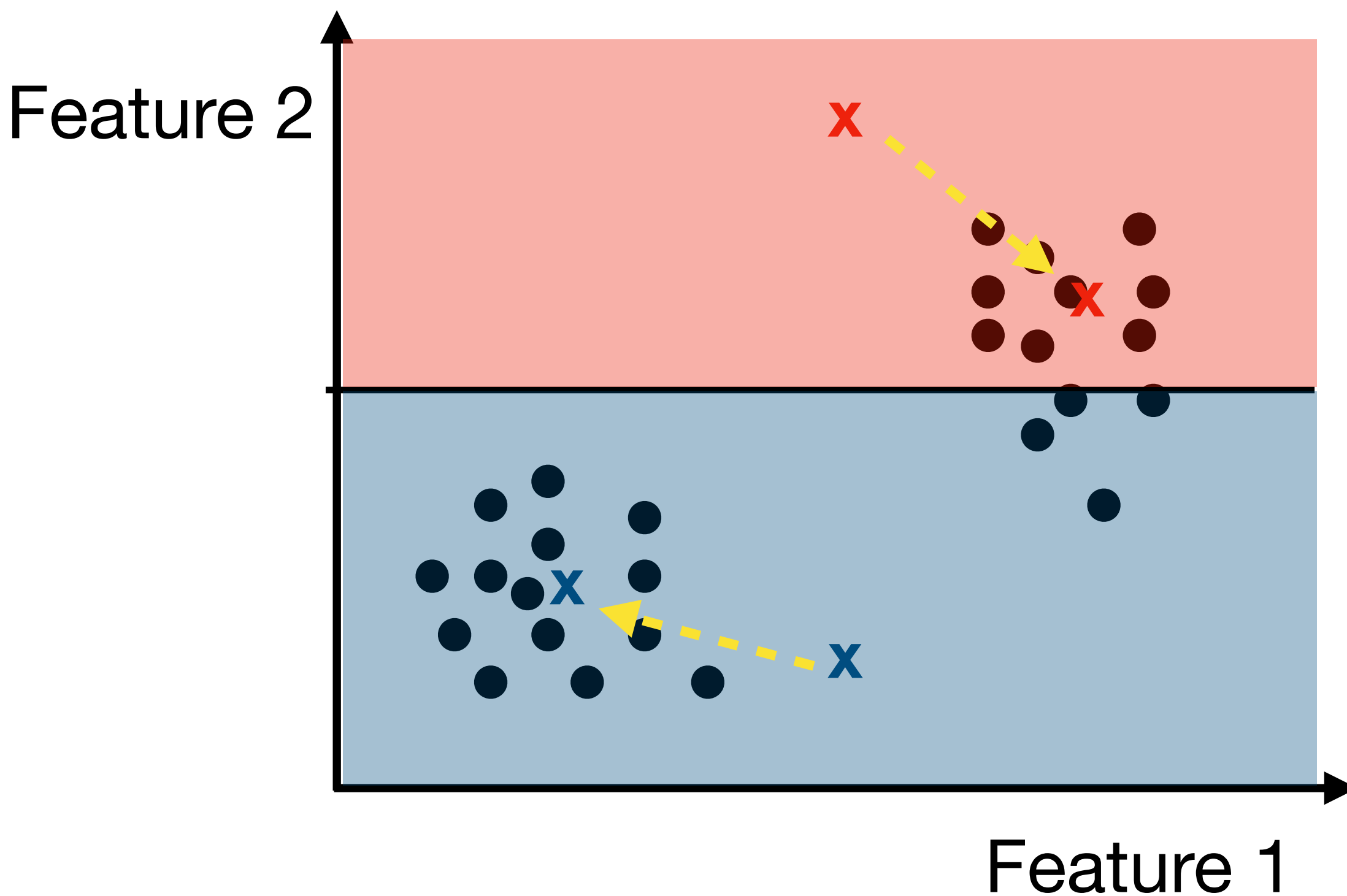
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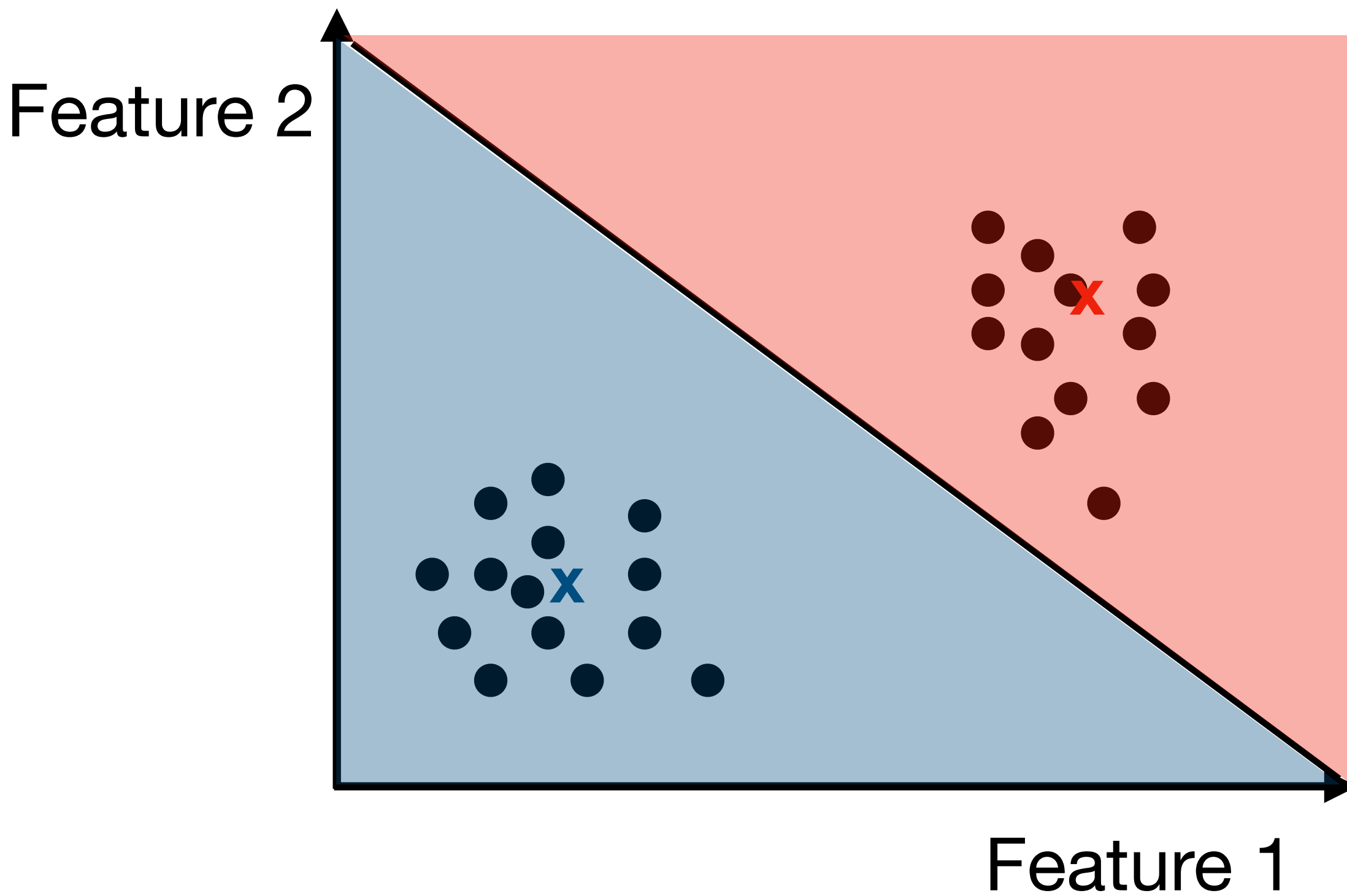
- Groups together data points with similar characteristics
- The methods for doing this typically involve distances in the plane of your observables



- One hyperparameter to be tuned: the number of clusters, k .
- Start with k centroids of clusters which are randomly initialised in the plane of the features
- Assign each point to a cluster based on the minimum distance to the centroid



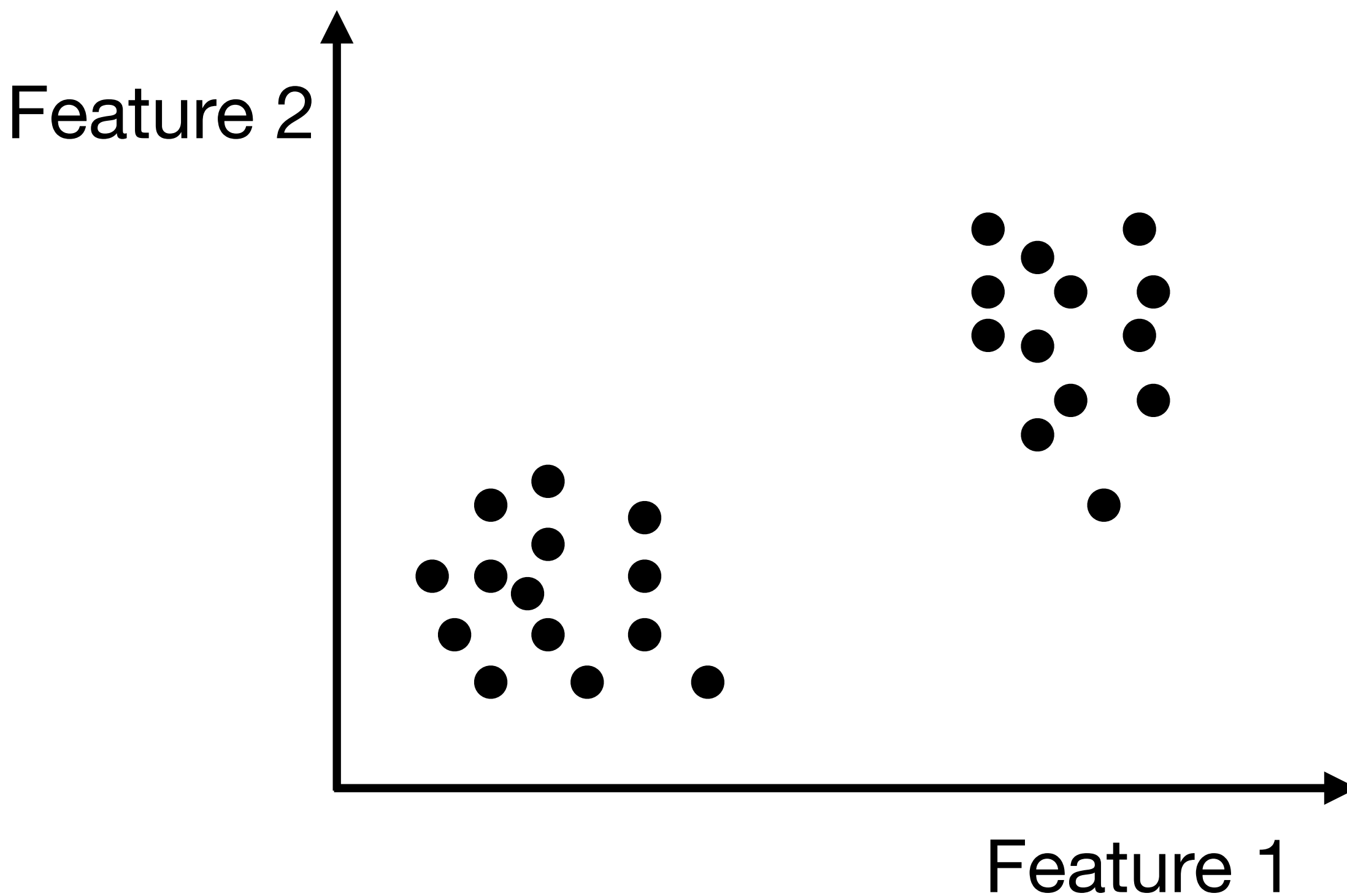
- Calculate the average of the points in the cluster region and make this average the new centroid
- Adjust the decision boundary accordingly by calculating the distances to the new centroids of every point



- We judge the algorithm to be converged to the number of clusters after the centroid changing after each iteration falls below a certain threshold

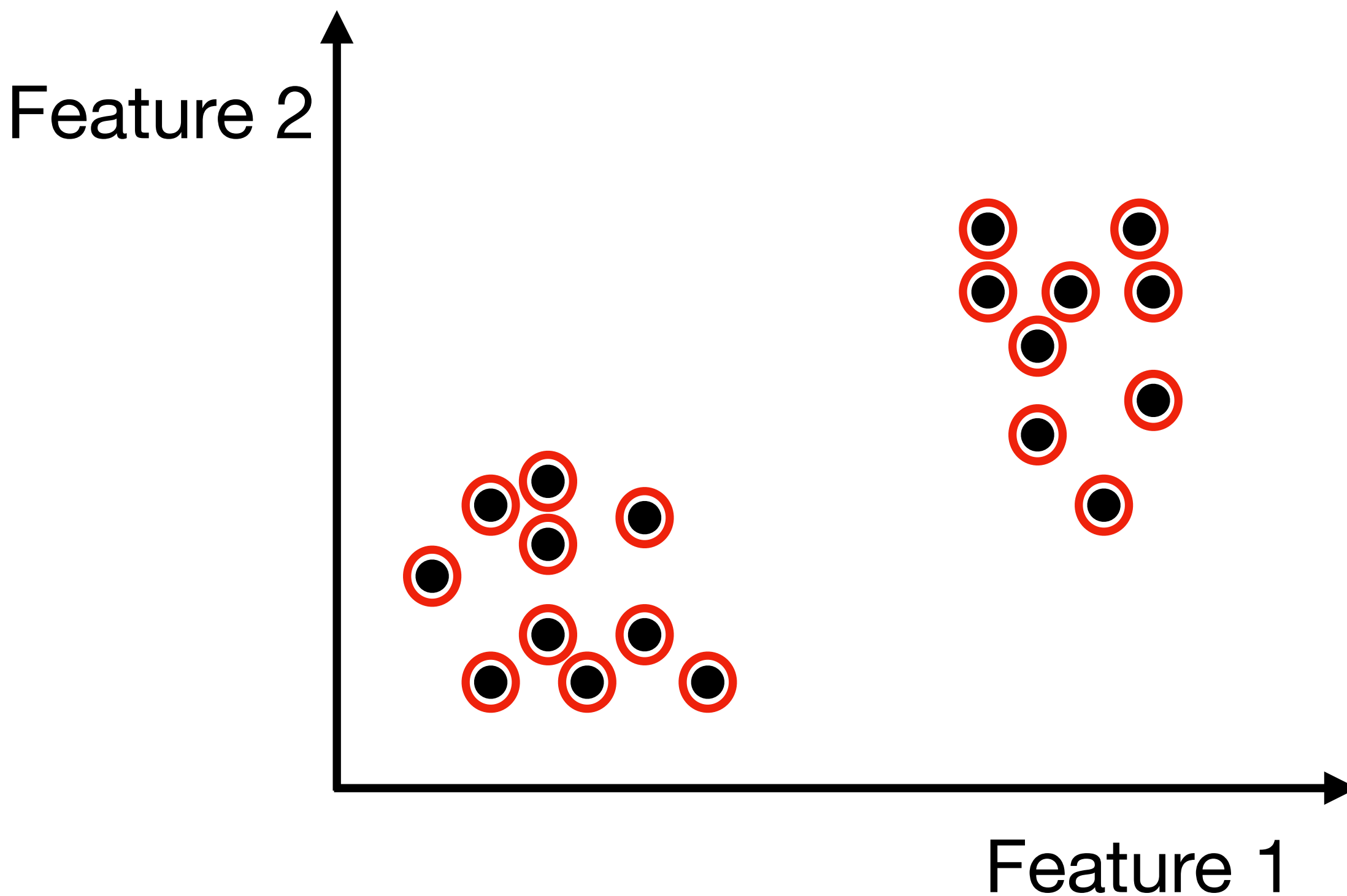
The method is simple:

- Each point starts as its own cluster
- The distances between the clusters are calculated
- The two closest clusters are merged into one cluster
- The process is repeated until all the points are in one large cluster

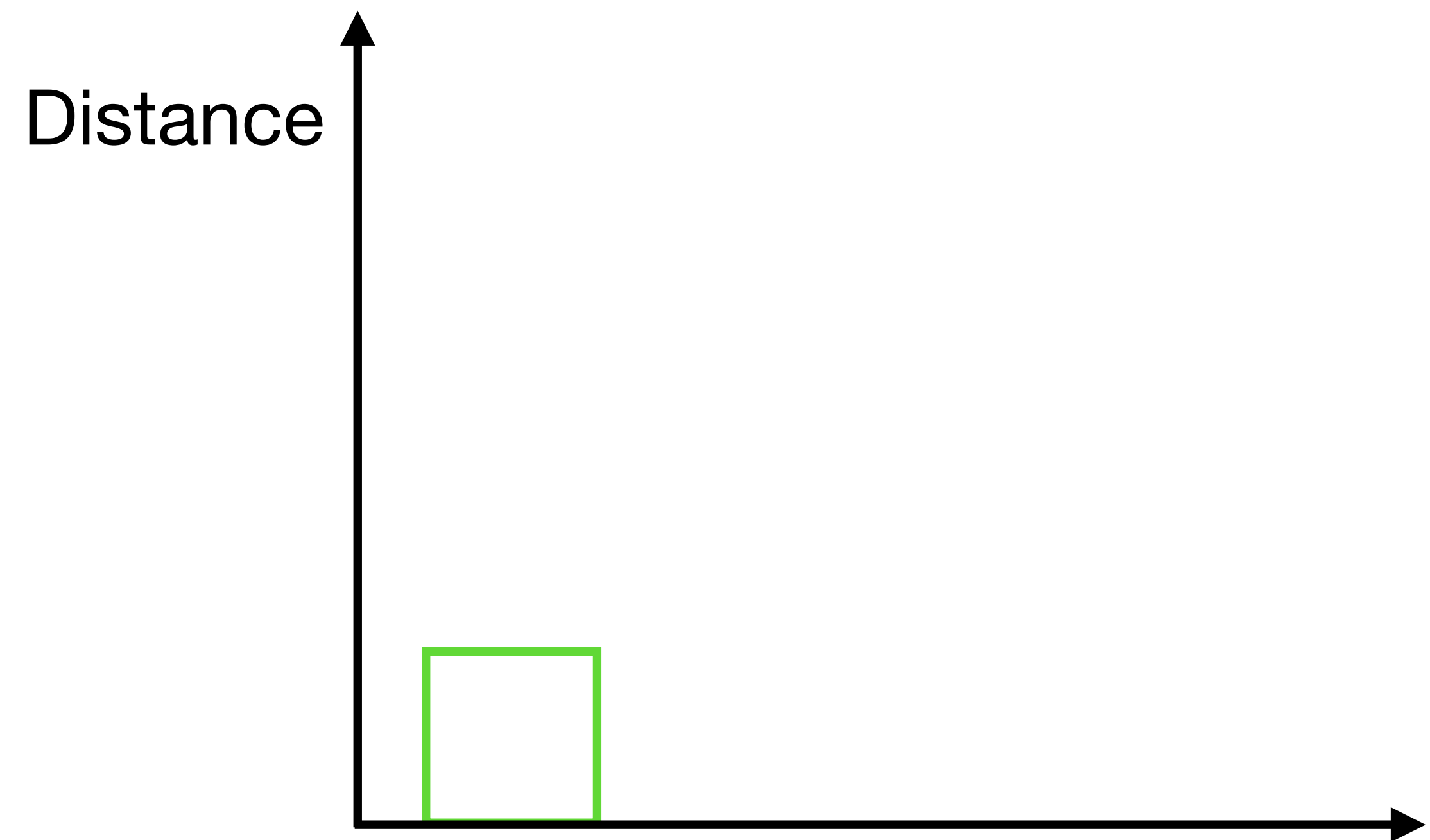
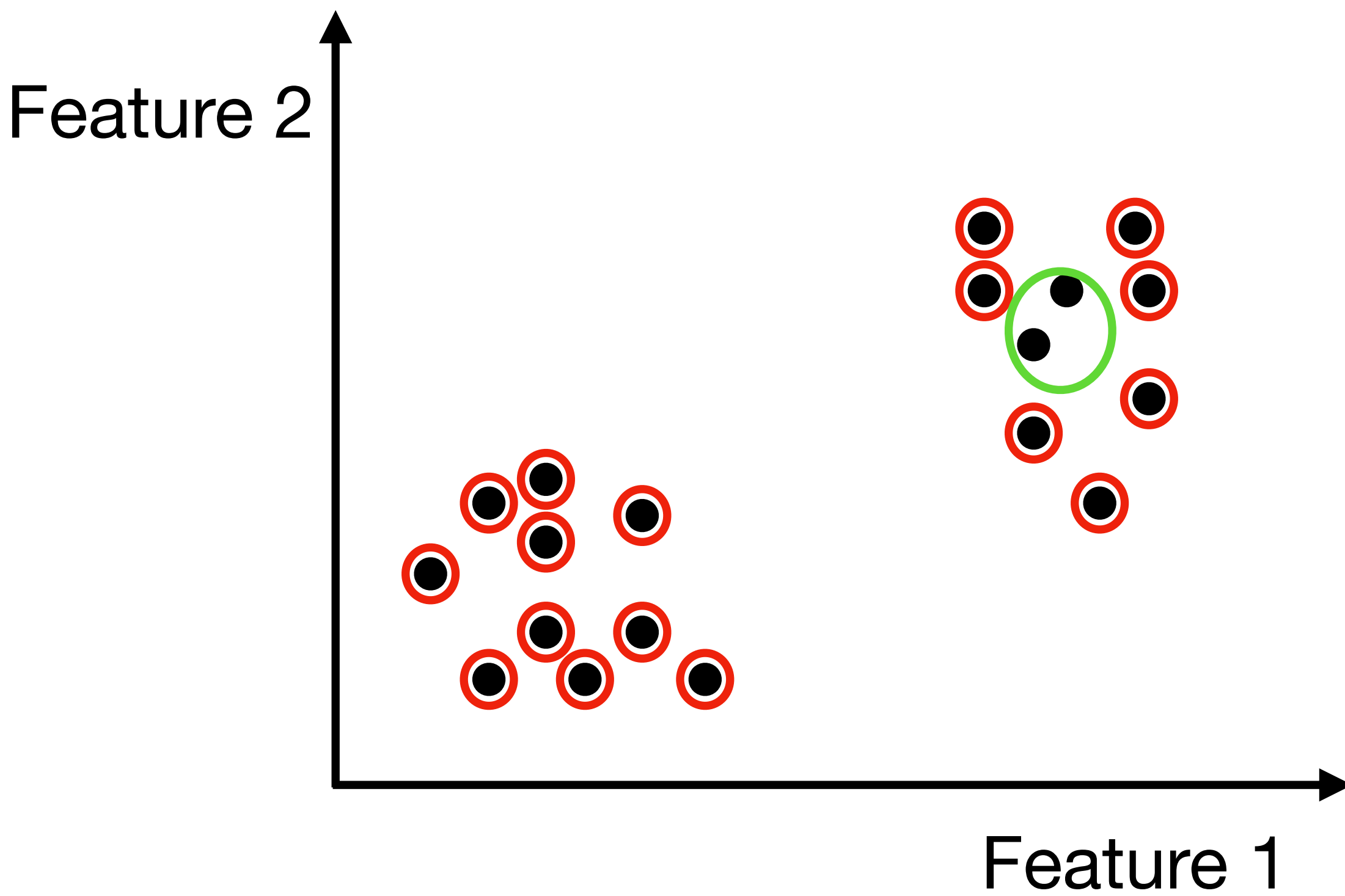


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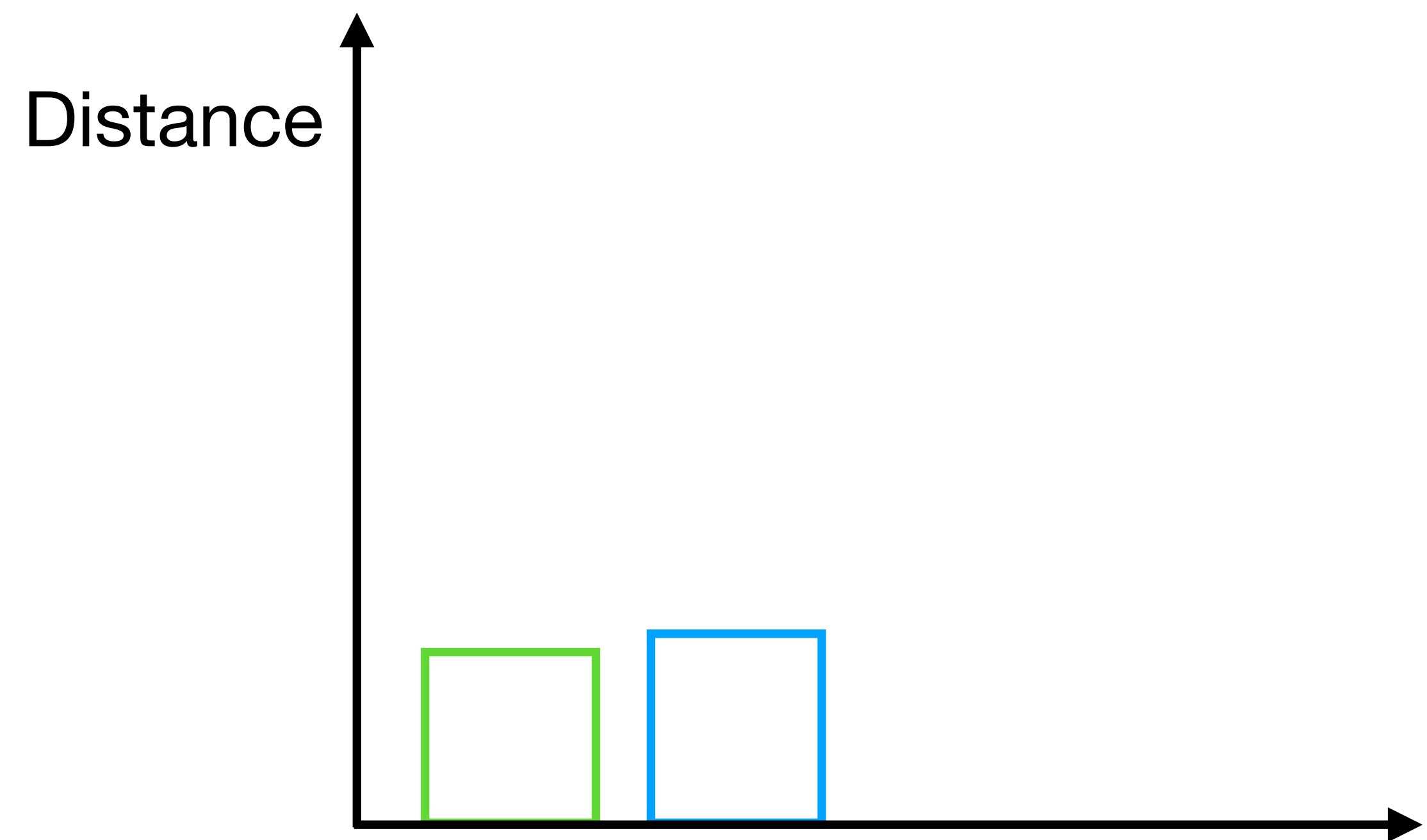
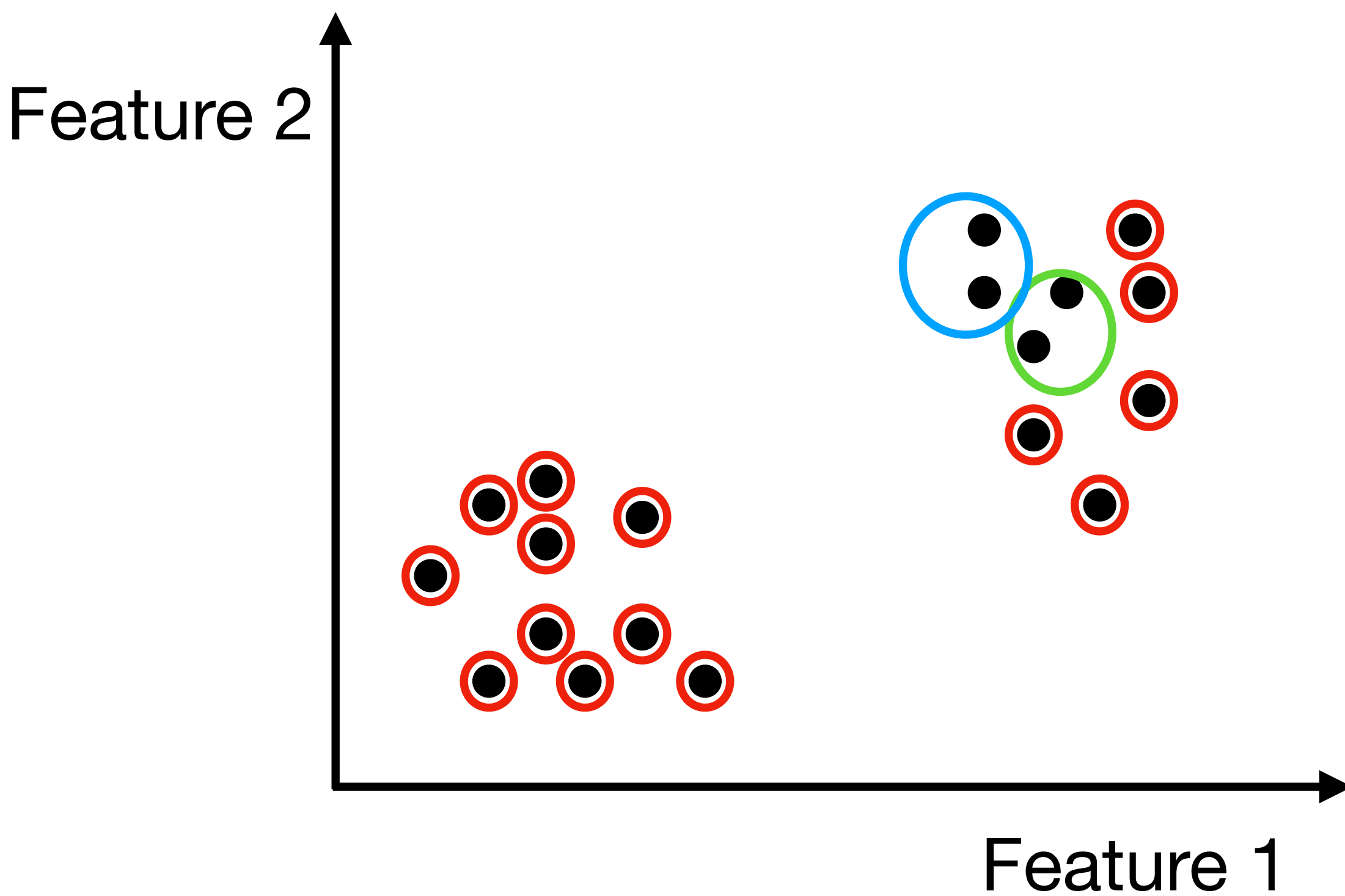
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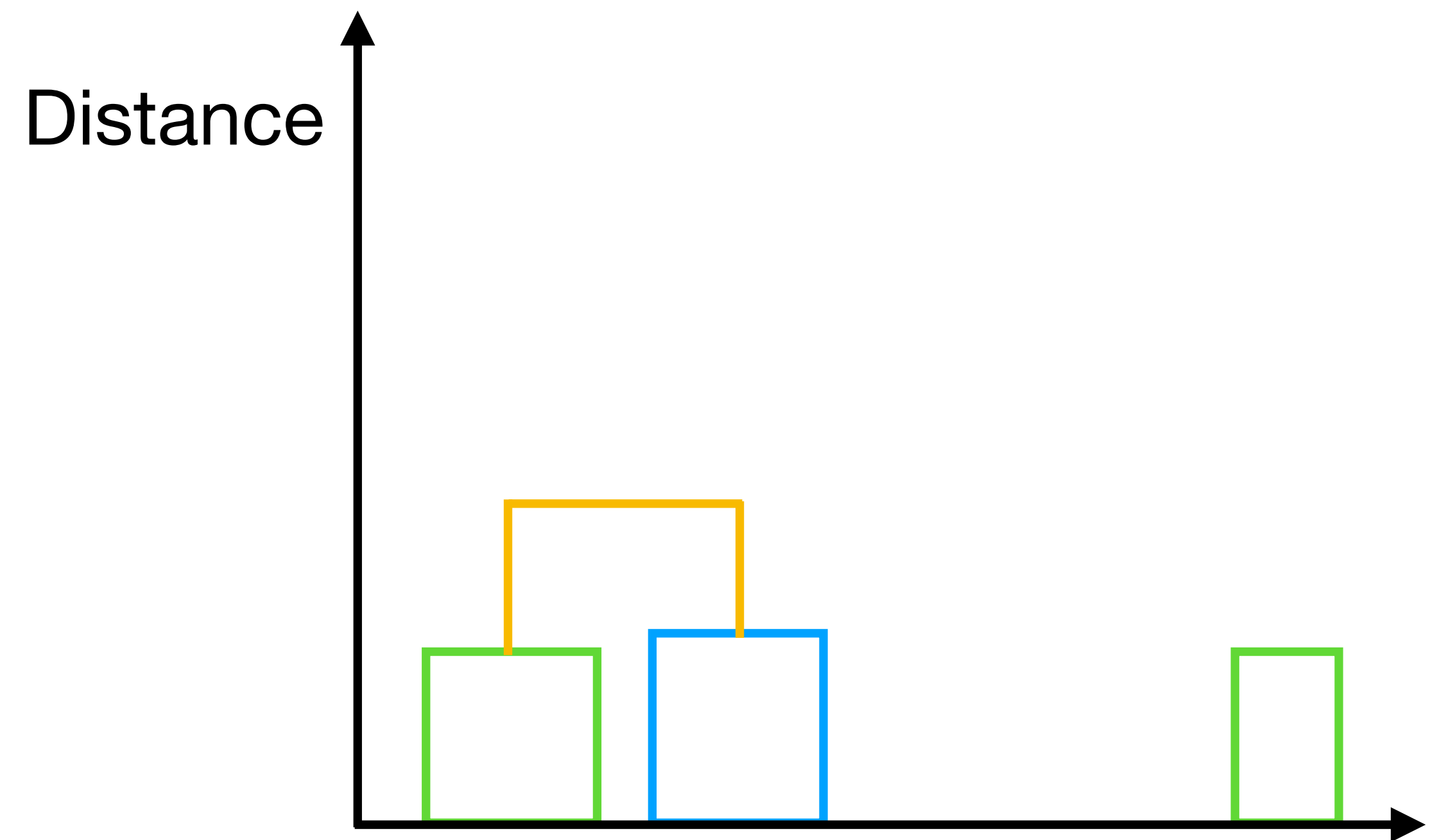
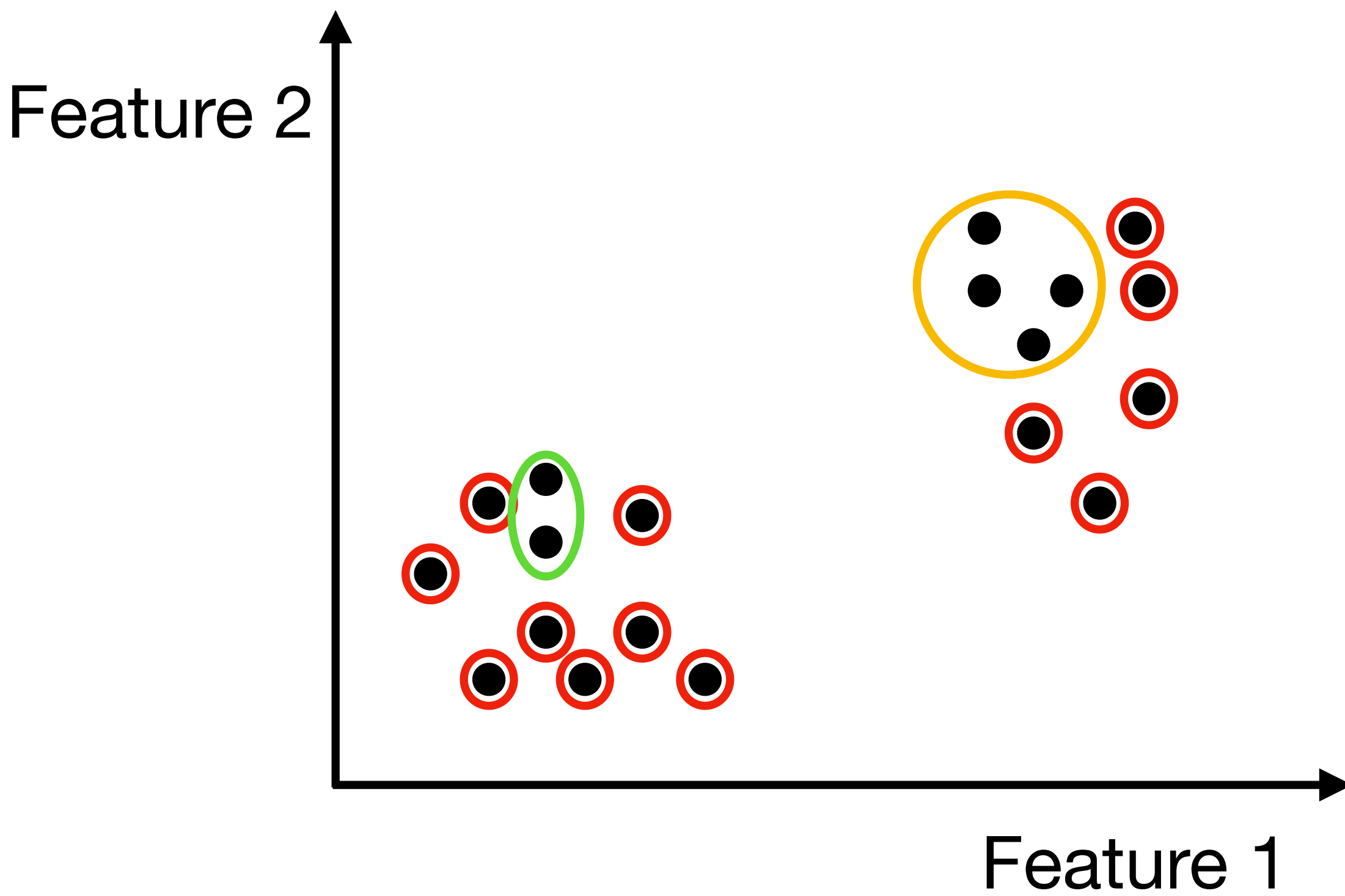
Dendrogram



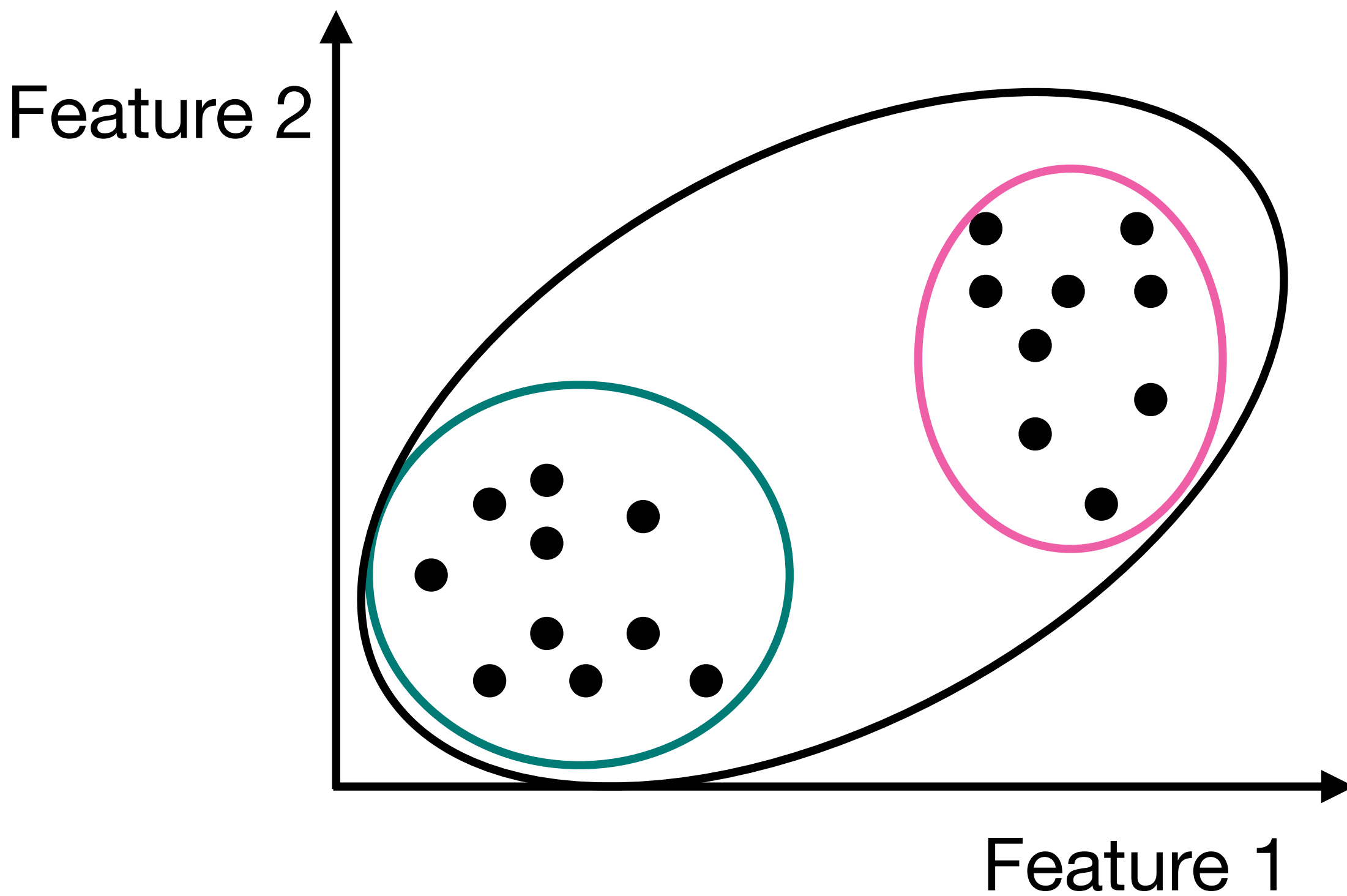
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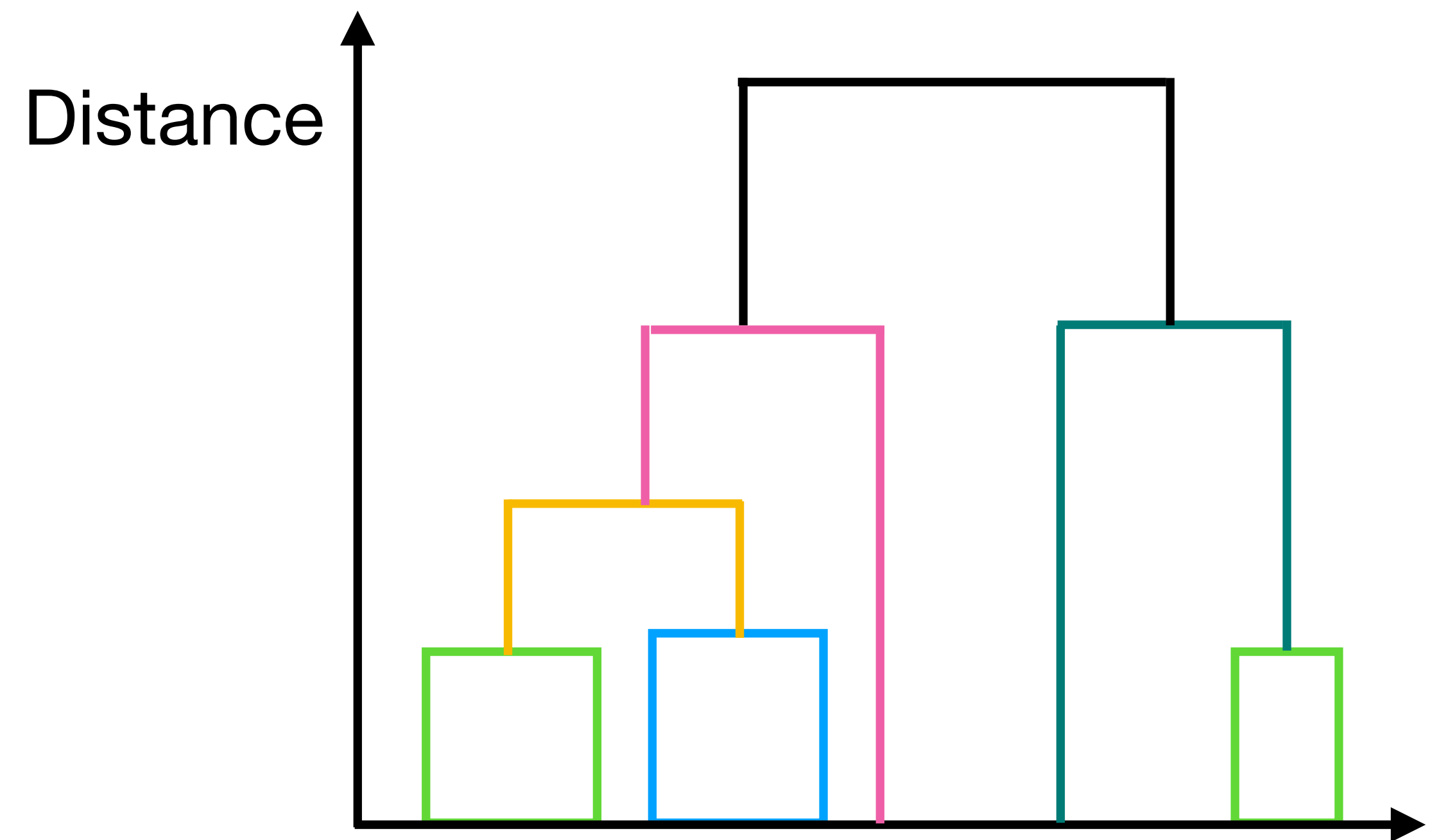
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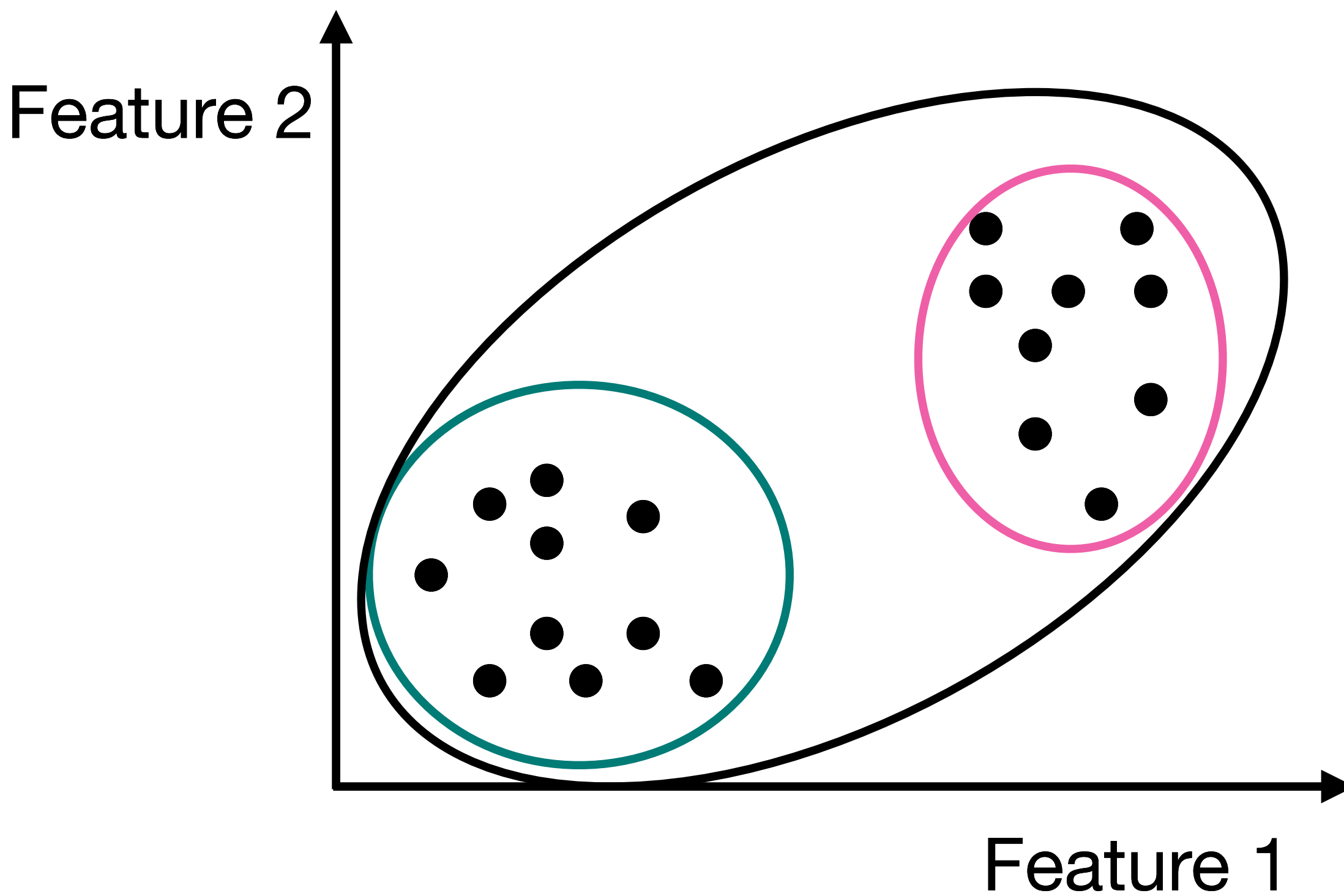


etc....



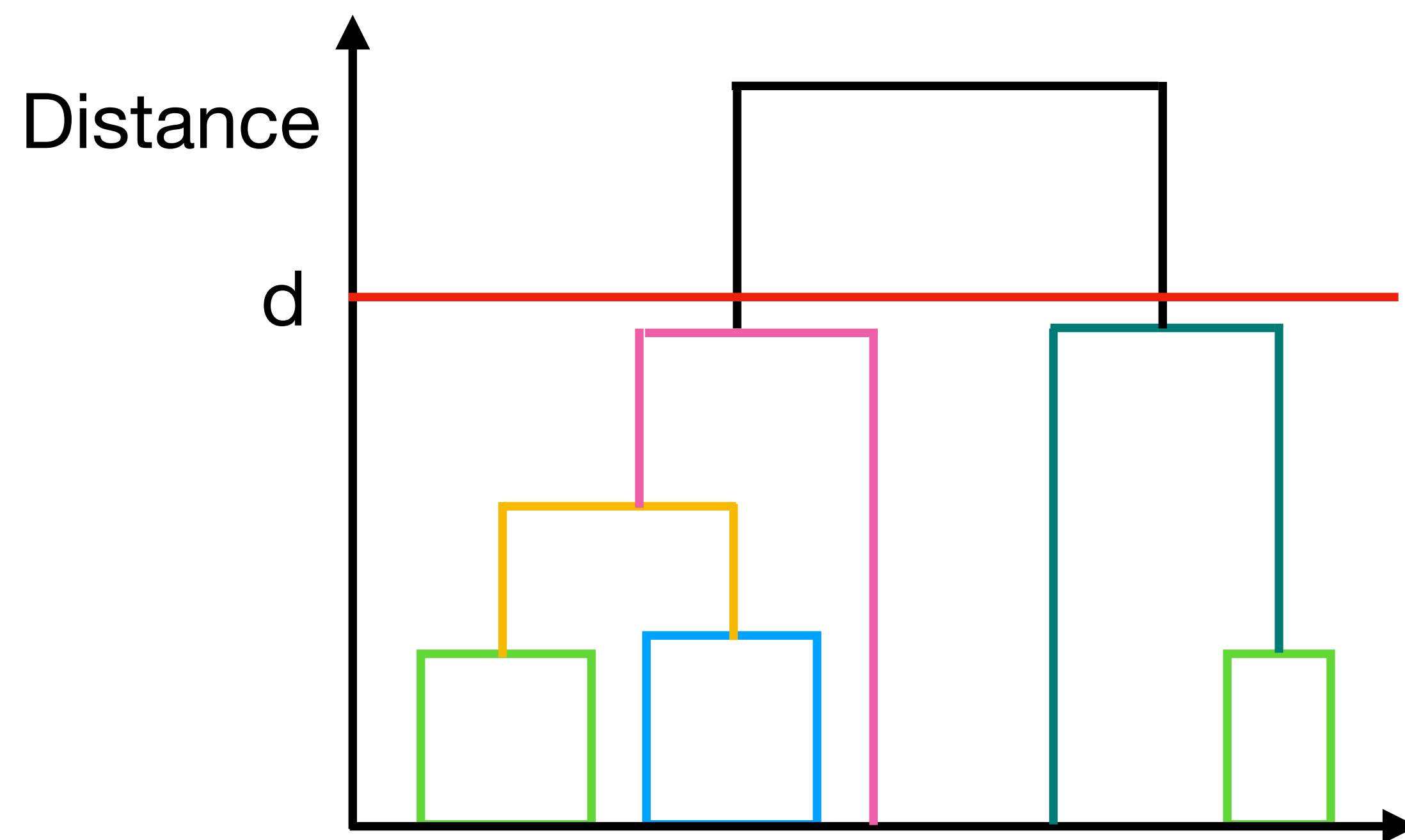
Dendrogram





There are 3 hyperparameters:

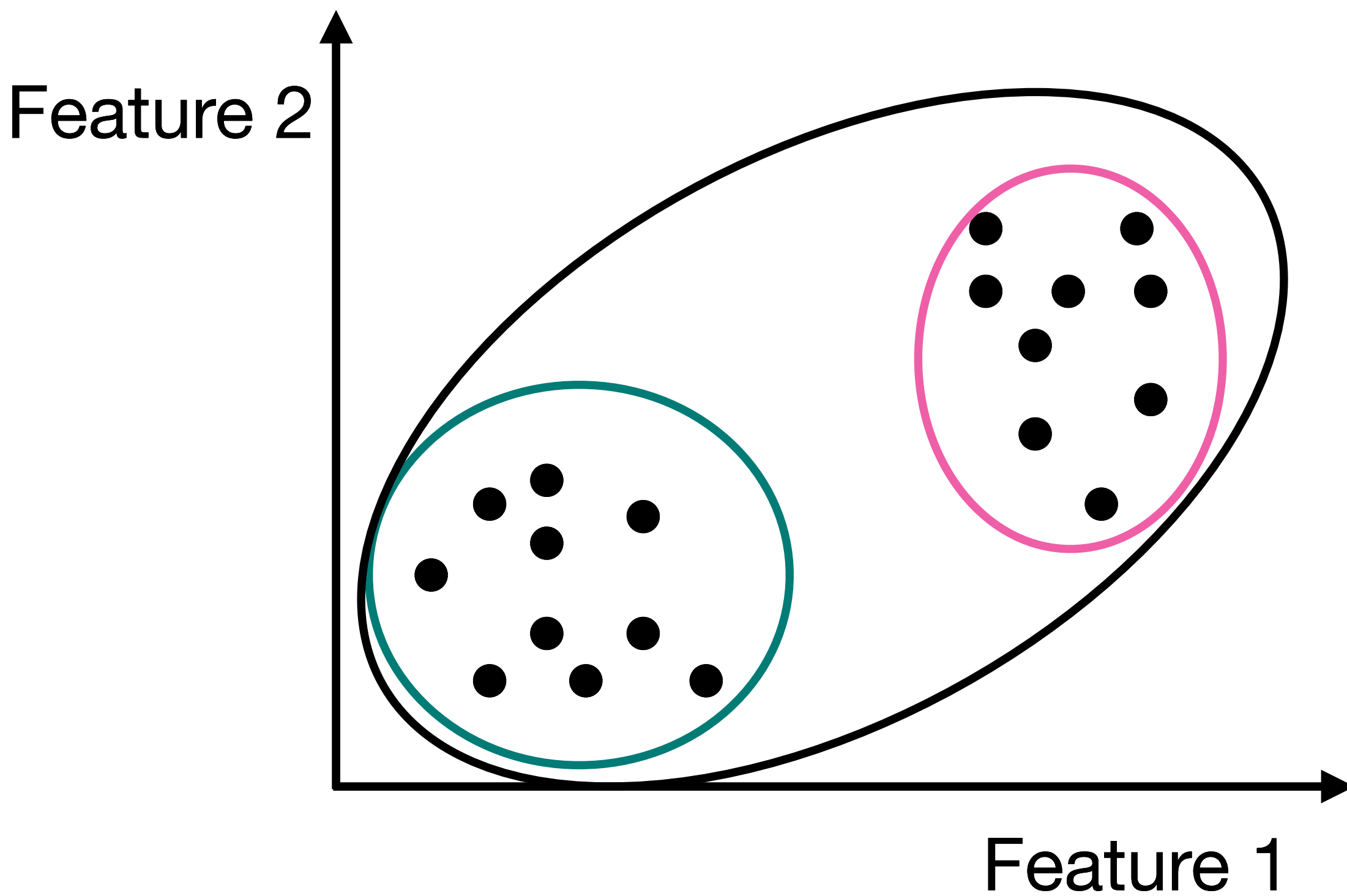
1. The number of clusters, k
e.g. in our example, it is obvious that $k = 2$,
however in reality, it may be more difficult and
you may need to experiment with k
Tuning k can be done by changing the cutoff
distance d in your dendrogram



There are 3 hyperparameters:

2. Cutoff distance, d

This is the distance at which you decide there is too large a distance between clusters for them to be part of one larger cluster



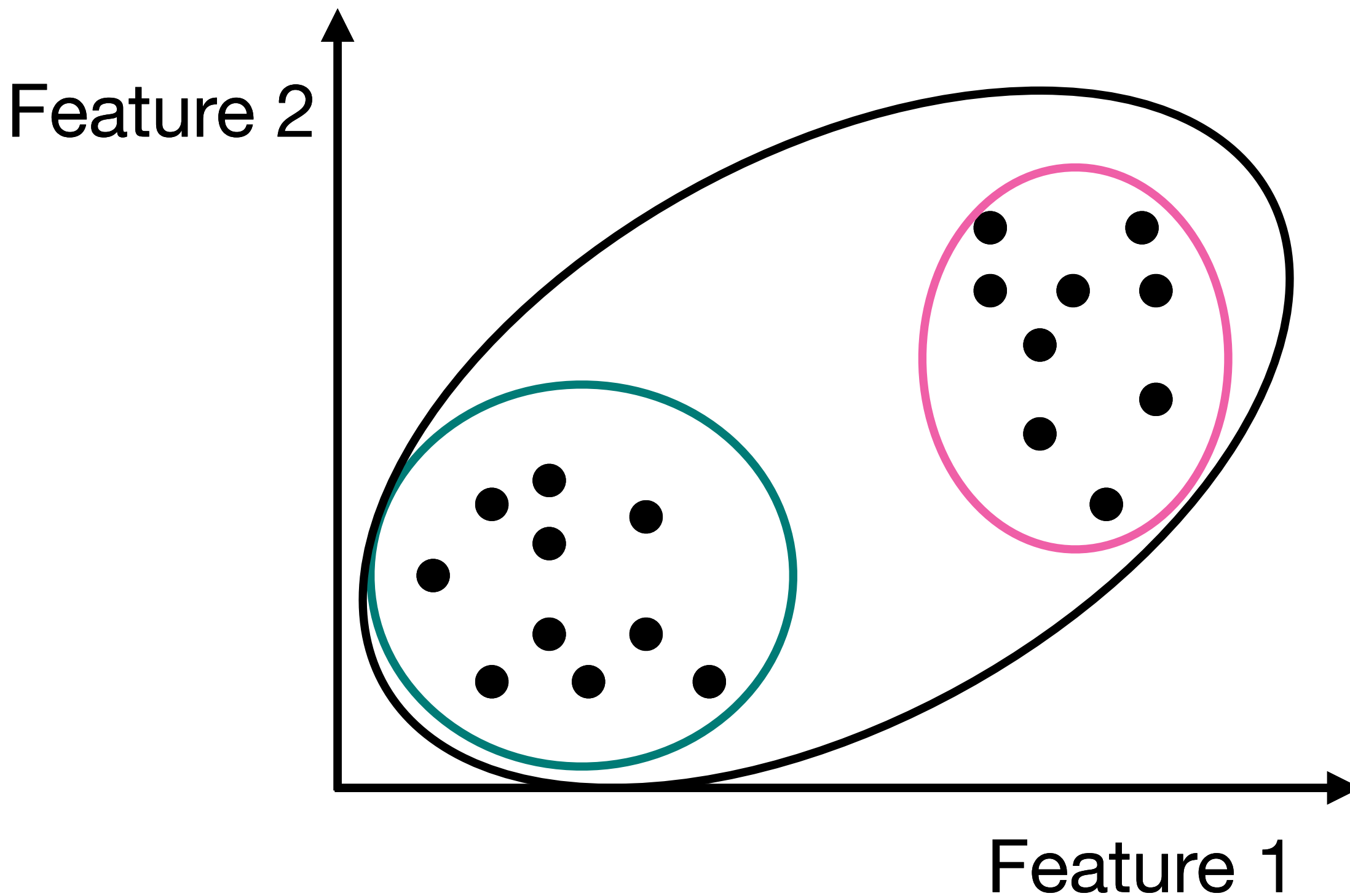
There are 3 hyperparameters:

3. The distance metric to calculate the distance between clusters e.g. Euclidean

$$d(\vec{x}_1, \vec{x}_2) = ||x_2 - x_1|| = \sqrt{\sum_{i=1}^n (x_1^i - x_2^i)^2}$$

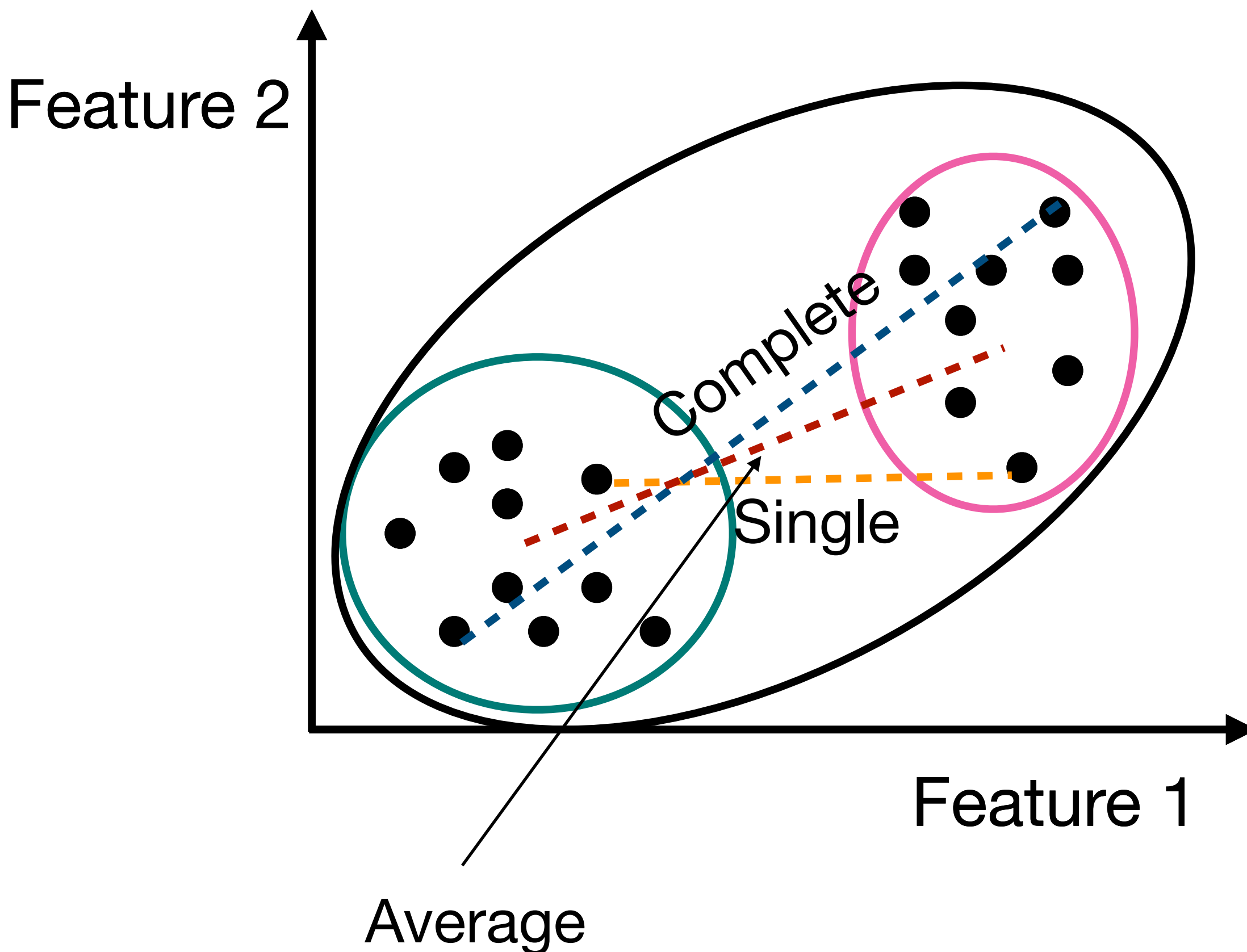
Linkage method:

- Finding distances between clusters is something also to be tuned (find which works best for you)
- When clusters have multiple points, defining the “distance” between two clusters is non-trivial
- How we do this is known as our algorithm’s **linkage method**



Examples of linkage methods:

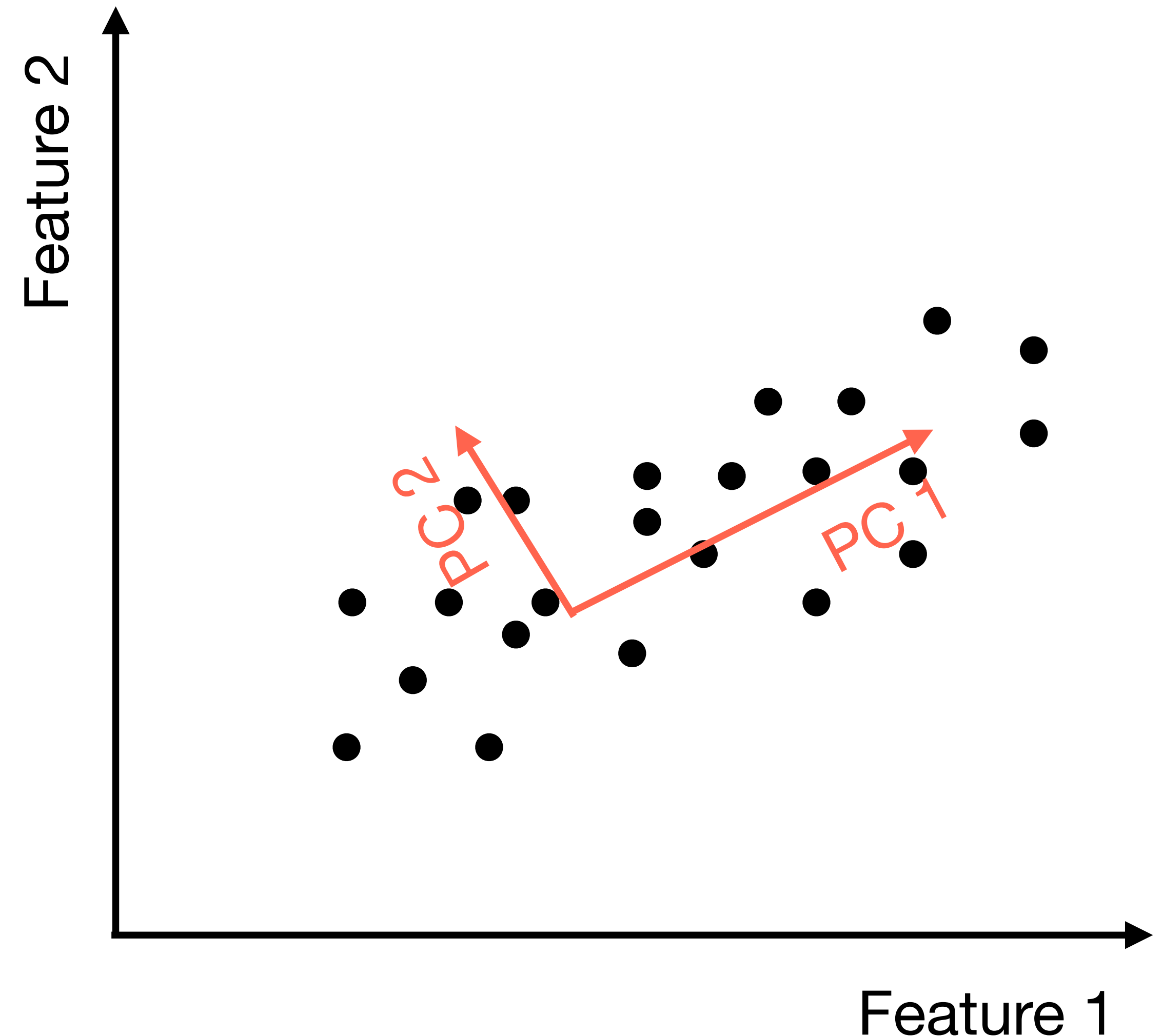
- Single — distance between two closest points
- Average — distance between average cluster position
- Complete — distance between two furthest points
- Ward — keeps growth in sum of squares as small as possible (requires Euclidean metric)



- Decomposition of data into important features and embedding a high-dimensional dataset into a lower-dimensional space.
- Important as certain features of your data may be redundant
- Improves performance of supervised learning
- Data compression/uncover complex trends

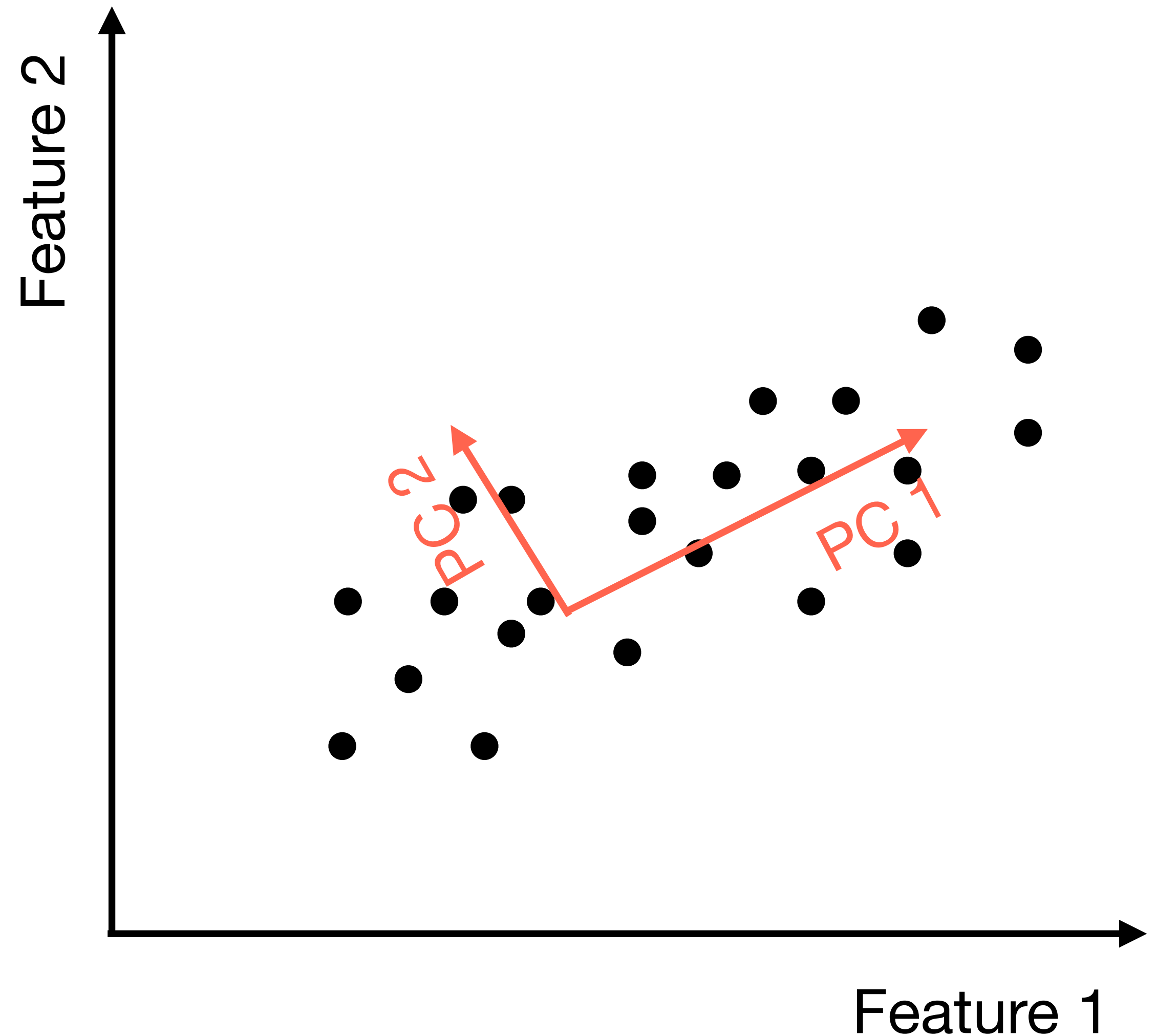
Principal Component Analysis (PCA)

- Orthogonal linear transformation to basis of so-called “principal components”
- These are the directions in the data with the most variance



Principal Component Analysis (PCA)

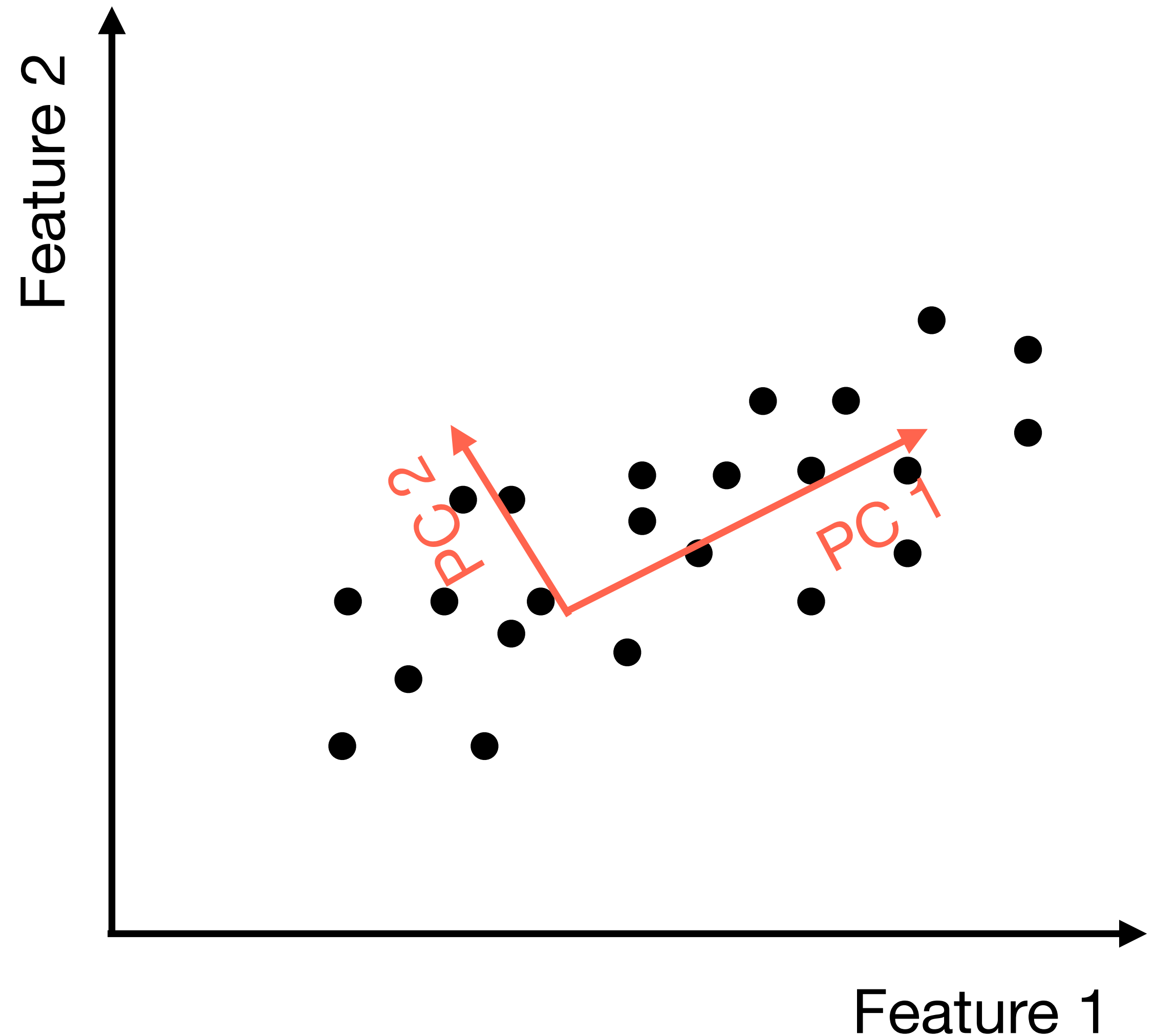
- The first principal component has the largest variance in the data
- Then each proceeding component has the largest variance of the remaining directions



Principal Component Analysis (PCA)

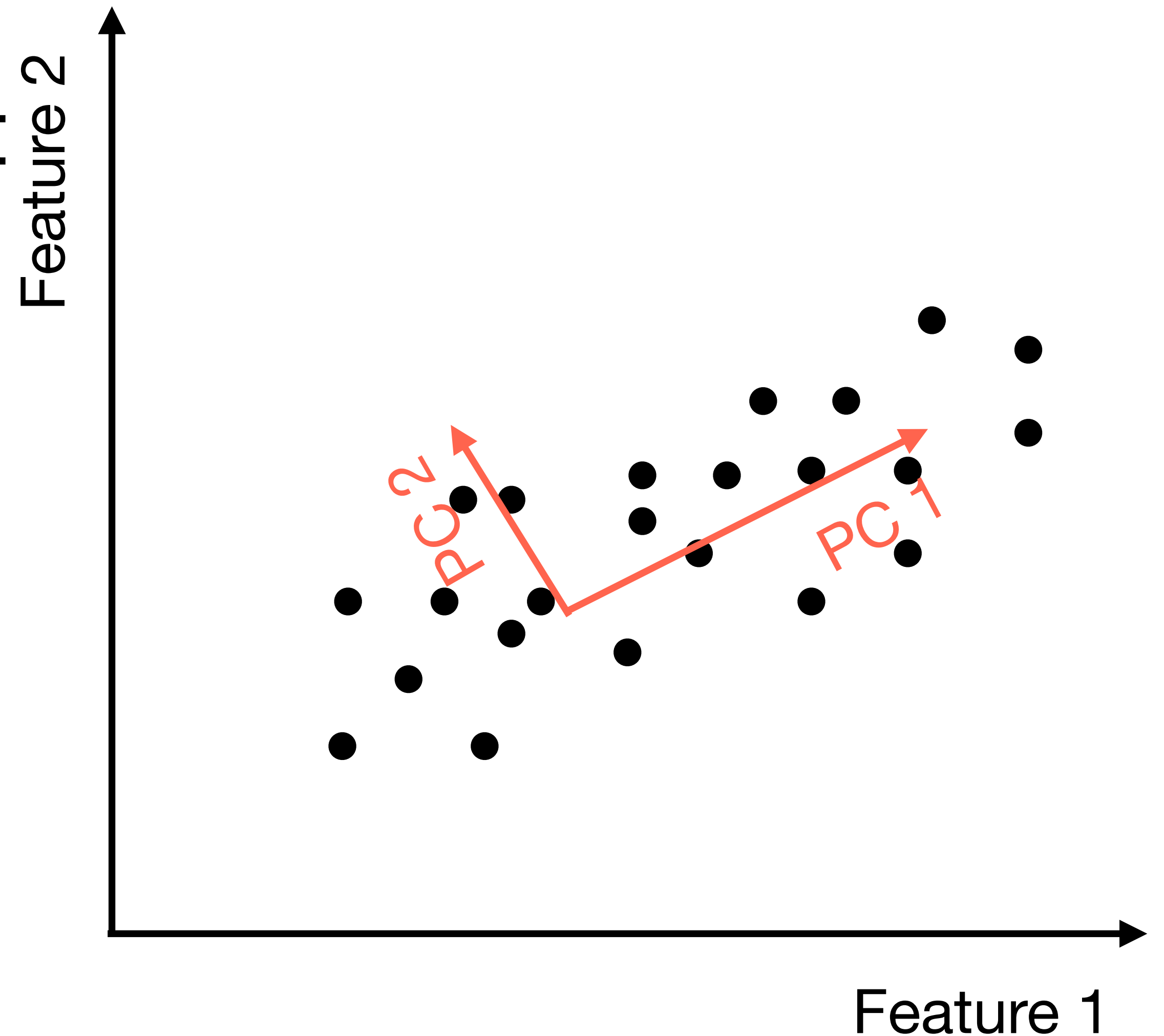
- PCA allows you to represent the data as a linear combination of the principal components

$$x = \sum_{i=1}^N A_i \times PC_i$$



Principal Component Analysis (PCA)

- Data can then be compressed by expressing it in terms of the most important principal components
- This also gives a low-dimensional representation



There are many other methods

- I have only explained two of a variety of techniques that can be used
- There are many other clustering and dimensionality reduction techniques to be used that may be more useful to your research
- Classical unsupervised learning can also be used for outlier detection and classification which we have not covered.
- <https://scikit-learn.org/> has good examples for many methods and an easy-to-use API
- sklearn also has a great cheatsheet: <http://blog.kaggle.com/2015/04/15/scikit-learn-video-2-setting-up-python-for-machine-learning/> on deciding what algorithm to use

- The exercises can be found on my GitHub: <https://bit.ly/2DtXei2> and are (hopefully) self-explanatory
- Feel free to email me with any questions about machine learning and/or the CDT (whether it be structure or courses or events):
j.armstrong.2@research.gla.ac.uk
- Honest feedback is appreciated!